# Progress Towards the First Measurement of Direct CP-Violation in $\mathbf{K} \to \pi\pi$ Decays From First Principles

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#### **RBC & UKQCD Collaboration**

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#### Outline

- Introduction
- $K \to (\pi \pi)_{I=2}$  calculation.
- $K \to (\pi\pi)_{I=0}$  calculation.
- Conclusions and Outlook

#### CP-Violation in the Standard Model

• Standard Model allows violation of CP via complex phase  $\delta$  in the CKM matrix.

$$\begin{pmatrix} \mathbf{d}' \\ \mathbf{s}' \\ \mathbf{b}' \end{pmatrix} = \begin{bmatrix} \mathbf{c}_{12} \mathbf{c}_{13} & \mathbf{s}_{13} \mathbf{e}^{i\delta} \\ -\mathbf{s}_{12} \mathbf{c}_{23} - \mathbf{c}_{12} \mathbf{s}_{23} \mathbf{s}_{13} \mathbf{e}^{i\delta} & \mathbf{c}_{12} \mathbf{c}_{23} - \mathbf{s}_{12} \mathbf{s}_{23} \mathbf{s}_{13} \mathbf{e}^{i\delta} & \mathbf{s}_{23} \mathbf{c}_{13} \\ \mathbf{s}_{12} \mathbf{s}_{23} - \mathbf{c}_{12} \mathbf{c}_{23} \mathbf{s}_{13} \mathbf{e}^{i\delta} & -\mathbf{c}_{12} \mathbf{s}_{23} - \mathbf{s}_{12} \mathbf{c}_{23} \mathbf{s}_{13} \mathbf{e}^{i\delta} & \mathbf{c}_{23} \mathbf{c}_{13} \end{bmatrix} \begin{pmatrix} \mathbf{d} \\ \mathbf{s} \\ \mathbf{b} \end{pmatrix}$$

- Manifests in 2 ways: Direct and Indirect
- Indirect CPV arises because weak eigenstates  $\neq$  CP eigenstates: e.g.  $K_S \propto K_1 + \bar{\epsilon} K_2$  where  $K_1$  and  $K_2$  are CP-even and CP-odd resp.
- Also direct CPV in decays of CP eigenstates:

$$K_1 ext{ (CP - even)} o \pi \pi \pi ext{ (CP - odd)}$$
  
 $K_2 ext{ (CP - odd)} o \pi \pi ext{ (CP - even)}$ 

# Brief interlude: lattice methods

- Discretize QCD Lagrangian in Euclidean space on finite volume.
- Integrate fermions out of path integral:

$$Z = \int dU \det(D[U]) \exp(-S_g[U])$$

- U are gauge links:  $U_{\mu} = e^{iaA_{\mu}^{a}T^{a}} \in SU(3)$
- Sample configurations of links from probability distribution Z using Monte Carlo methods.

#### Lattice measurements

 Measure amplitudes on each link configuration and average.

$$\int d^3 \vec{x} \, \langle 0 | \bar{d}(x) \gamma^5 u(x) \bar{u}(0) \gamma^5 d(0) | 0 \rangle$$

$$= \frac{1}{N} \sum_{i=1}^{N} \int d^3 \vec{x} \, \text{tr} \left( \gamma^5 D_d^{-1}(0, x) [U_i] \gamma^5 D_u^{-1}(x, 0) [U_i] \right)$$

$$= a_0 e^{-m_\pi x_4} + a_1 e^{-E_1 x_4} + \dots$$

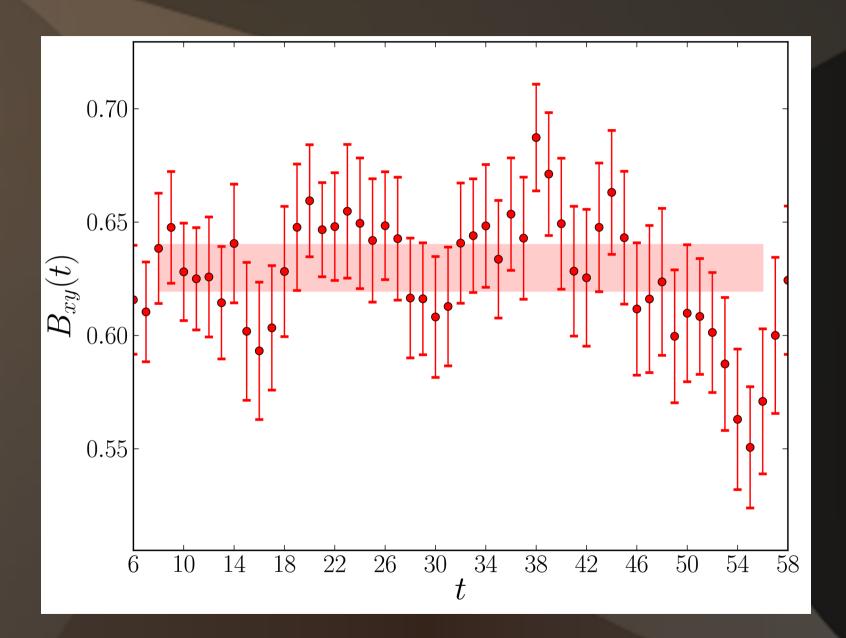
- Ground state of system extracted in limit of large time separation.
- Excited state with energy  $E_i$  (i > 0) requires multiexponential fits to time dependence – typically very noisy and should be avoided if possible!

### Indirect CP-Violation on the Lattice

- Indirect CPV measure  $\epsilon$  determined accurately from experiments:  $\epsilon = \frac{A(K_L \to (\pi\pi)_{I=0})}{A(K_S \to (\pi\pi)_{I=0})}$
- Theoretically  $\epsilon \propto G_F^2 M_W^2 B_K(\mu) S(\mu)$  where  $S(\mu)$  are perturbative Wilson coefficients and  $B_K(\mu)$  contains the non-perturbative QCD contribution.
- Both factors are renormalization scheme dependent but their product is scheme invariant.
- On lattice we can measure  $B_K$  through

$$B_K \propto \langle \bar{K}^0 | \mathcal{H}_{\text{eff}}(\Delta S = 2) | K^0 \rangle$$

Modern calculations at %-scale accuracy.



# $\mathbf{K} ightarrow \pi\pi$ Decays

- Direct CP-violation first observed in  $K \to \pi\pi$  decays.
- Two types of decay:

$$\begin{array}{cccc} \Delta I = 3/2 & :K^+ & \rightarrow (\pi^+\pi^0)_{I=2} & \text{with amplitude } A_2 \\ \Delta I = 1/2 & :K^0 & \rightarrow (\pi^+\pi^-)_{I=0} \\ & K^0 & \rightarrow (\pi^0\pi^0)_{I=0} & \text{with amplitude } A_0 \end{array}$$

- Direct CP-violation:  $\epsilon'=rac{i\omega e^{i(\delta_2-\delta_0)}}{\sqrt{2}}\left(rac{{
  m Im}A_2}{{
  m Re}A_2}-rac{{
  m Im}A_0}{{
  m Re}A_0}
  ight)$  where
  - $\omega = \mathrm{Re}A_2/\mathrm{Re}A_0$  and  $\delta_I$  are strong scattering phase shifts
- $\epsilon'$  is highly sensitive to BSM sources of CPV.
- Strong interactions very important origin of the so-called  $\Delta I=1/2$  rule: preference to decay to I=0 final state.

#### $\mathbf{K} o \pi\pi$ on the lattice

 Multi-particle states in a finite box very different from infinite-volume states:

$$|\pi\pi\rangle_{\text{latt}} = c_0|\pi\pi \ (l=0)\rangle_{\text{phys}} + c_4|\pi\pi \ (l=4)\rangle_{\text{phys}} + \dots$$

- Until recently not known how to relate lattice amplitude to physical amplitude. [Lellouch&Luscher]
- Energy spectrum is volume-dependent; need large physical volume for realistic kinematics.
- Also need small lattice spacing to avoid large discretization errors.
- Large volume + small lattice spacing = expensive!
- Only recently become viable.

 $K \rightarrow (\pi\pi)_{I=2}$  Calculation

#### Lattice Determination

• As with  $B_K$ , amplitude  $A_2$  is combination of renormalization-scheme dependent perturbative Wilson coeffs  $C_i(\mu)$  and non-perturbative matrix elements  $M_i(\mu)$ :

$$A_2 \propto G_F V_{ud} V_{us} \sum_{i=1}^{10} C_i(\mu) M_i(\mu)$$

- $M_i = \langle (\pi^+ \pi^0)_{I=2} | Q_i | K^0 \rangle^{i=1}$
- $Q_i$  are weak effective four-quark operators.
- Renormalization performed non-perturbatively in intermediate regularization-independent momentum scheme (RI-MOM), matched to  $\overline{\rm MS}$  at high energies to avoid perturbative truncation errors.

# Achieving Physical Kinematics

- $m_{\pi} = 135 \text{ MeV}$  and  $m_{K} = 500 \text{ MeV}$ : need moving pions in final state to conserve energy.
- Ground state of  $\pi\pi$  system has stationary pions.
- As previously mentioned, extracting excited states is very hard. Can we avoid this? Yes!

# Physical Kinematics

Instead impose antiperiodic BCs on d-quark propagator.
 Changes finite-volume momentum discretization:

$$p = \frac{2\pi n}{L} \to \frac{(2n+1)\pi}{L}$$

- Minimum d-quark momentum is  $\pi/L$ : charged pion ground state has momentum! But...
- For neutral pion the momenta can cancel, s.t. ground state is stationary. Desired state is  $\pi^+\pi^0$ , so this does not work. However....
- Wigner-Eckart theorem:

$$\langle (\pi^+\pi^0)_{I=2}|Q^{\Delta I_z=1/2}|K^+\rangle = \frac{\sqrt{3}}{2}\langle (\pi^+\pi^+)_{I=2}|Q^{\Delta I_z=3/2}|K^+\rangle$$
• APBCs on d-quark break isospin symmetry allowing mixing

• APBCs on d-quark break isospin symmetry allowing mixing between isospin states: however  $\pi^+\pi^+$  is the only charge-2 state hence it cannot mix.

#### Results

- RBC & UKQCD recently published (arXiv:1111.1699) calculation of  $\Delta I = 1/2$  decay using:
  - 2+1f domain wall fermions on a  $32^3 \times 64 \times 32$  lattice with  $a^{-1}=1.37(1)~{\rm GeV}.$
  - Near physical pions:  $m_\pi^{PQ} \sim 140 \; {
    m MeV}, m_\pi^{
    m uni} \sim 170 \; {
    m MeV}$
  - Energy conserving decays
- Determined

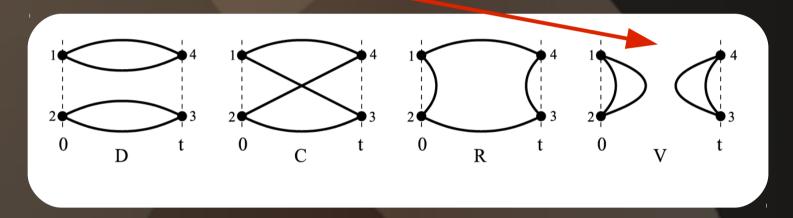
$$ReA_2 = [1.436(62)_{stat}(258)_{sys}] \times 10^{-8} \text{ GeV}$$
  
 $ImA_2 = -[6.83(51)_{stat}(1.30)_{sys}] \times 10^{-8} \text{ GeV}$ 

- Large systematic error of which 75% is discretization error: continuum limit needed.
- Currently generating multiple larger, finer lattices to get better control of this error.

 $K \rightarrow (\pi\pi)_{I=0}$  Calculation

# Challenges: part 1

- Measuring  $A_0$  is considerably more challenging.
- Measure both  $K^0 \to \pi^+\pi^-$  and  $K^0 \to \pi^0\pi^0$ .
- $\pi\pi$  state has vacuum quantum numbers, hence there are disconnected diagrams:



- Need large statistics and many source positions (or A2A/AMA propagators) to resolve.
- With Blue Gene/Q resources we can now perform such calculations with large enough physical volumes.

# Challenges: part 2

- For  $\Delta I = 1/2$  the Wigner-Eckart trick cannot be used.
- If we stay with APBC on d-quarks, isospin-breaking would allow mixing between I=0 and I=2 final states.
- I=0 state needs moving  $\pi^0$ , but momentum cancels in  $d\bar{d}$ .
- Need to apply BCs that commute with isospin and produce moving  $\pi^0$  as well as  $\pi^+$  and  $\pi^-$ .
- Can we conceive boundary conditions that satisfy these criteria? Yes: G-parity.

# G-Parity Boundary Conditions

• G-parity is a charge conjugation followed by a 180 degree isospin rotation about the y-axis:

Wiese, Nucl.Phys.B375, (1992)

$$\hat{G}=\hat{C}e^{i\pi\hat{I}_y}~:~\hat{G}|\pi^\pm\rangle=-|\pi^\pm\rangle \label{eq:Gamma} \begin{array}{c} \text{Kim, arXiv:hep-lat/0311003}\\ \hat{G}|\pi^0\rangle=-|\pi^0\rangle \end{array}$$

- Pions are all eigenstates with e-val -1, hence G-parity BCs make pion wavefunctions antiperiodic, with minimum momentum  $\pi/L$ .
- G-parity commutes with isospin

#### Kaons

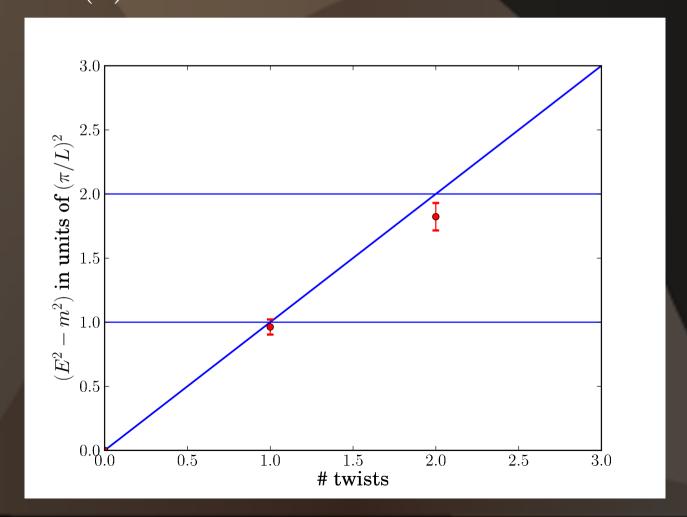
- $K \to \pi\pi$  calculation needs stationary  $K^0$ .
- Need an eigenstate with e-val +1 for periodic BCs and hence  $p_{\min} = 0$
- $\frac{1}{\sqrt{2}}(\bar{s}d + \bar{d}s)$  is not a G-parity eigenstate. Introduce 'strange isospin' (I'): s-quark in doublet  $\begin{pmatrix} s' \\ s \end{pmatrix}$
- A neutral kaon-like state:

$$K'_0 = \frac{1}{2}(\bar{s}d + \bar{d}s + \bar{s}'u + \bar{u}s')$$

is an eigenstate of 'modified G-parity':  $\hat{G} = \hat{C}e^{i\pi\hat{I}_y}e^{i\pi\hat{I}'_y}$  with eval + 1.

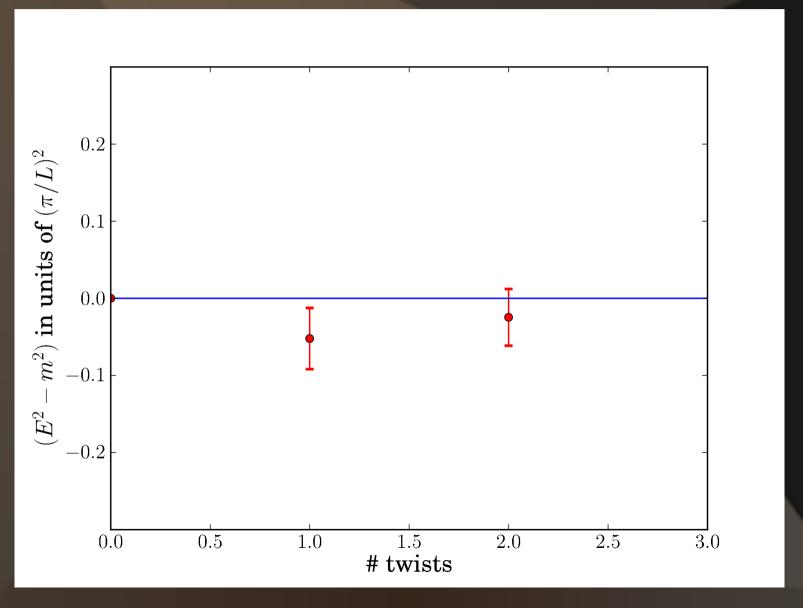
# Results: Pion Dispersion Relation

- Generated  $16^3 \times 32$  fully dynamical test ensembles with G-parity BCs in 0,1,2 directions.
- $a^{-1} = 1.73(3) \text{ GeV}$   $m_{\pi} \sim 420 \text{ MeV}$



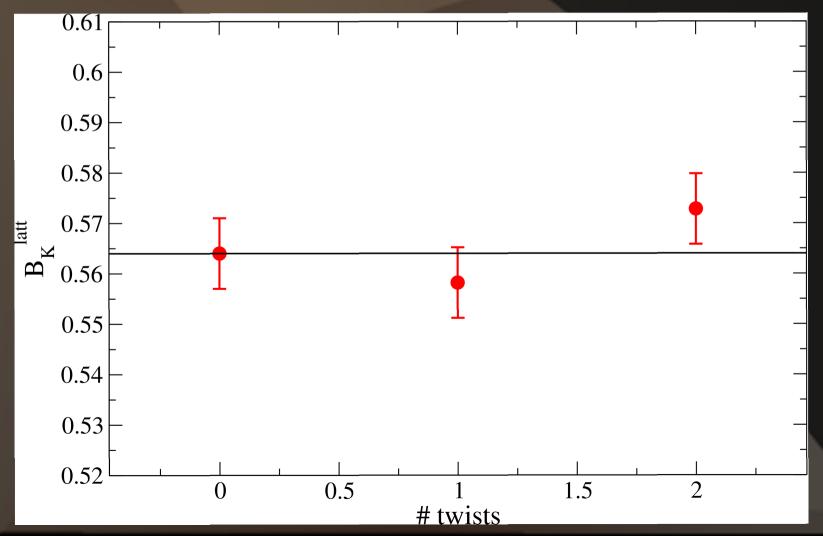
# Results: Kaon Dispersion Relation

Stationary kaon states demonstrated:



#### Results: $B_K$

•  $\bar{K}^0 \leftrightarrow K^0$  mixing amplitude shown to be independent as expected. These 4-quark effective vertices are similar to those used in  $K \to \pi\pi$  calculation, hence this is a valuable demonstration.



# Conclusions and Outlook

#### Conclusions and Outlook

- Lattice calculations have the potential to lead to great breakthroughs in our understanding of kaon phenomenology, in particular CP-violation.
- In the near future we will begin generating G-parity ensembles with large physical volumes and physical quark masses for a calculation of the  $\Delta I = 1/2$   $K \to \pi\pi$  amplitude.
- Combining with our existing measurement of the  $\Delta I = 3/2$  amplitude will give the first *ab initio* determination of  $\epsilon'$ . Could potentially lead to discovery of new BSM physics.

# Gauge Field Boundary Conditions

• d-field becomes  $C\bar{u}^T$  across the boundary. Consider a bilinear on the boundary under a gauge transformation :

$$\bar{d}(L-1)U_y(L-1)C\bar{u}^T(0)$$

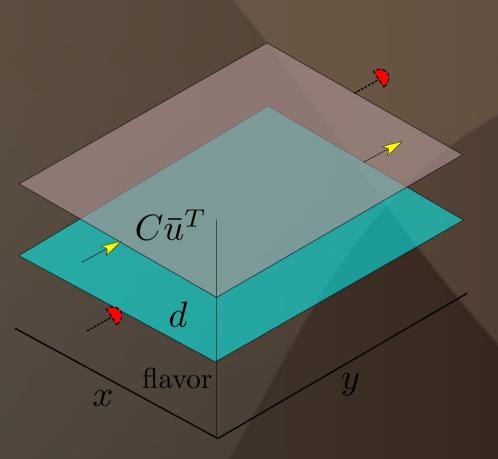
$$\longrightarrow \bar{d}(L-1)V^{\dagger}(L-1)U_y(L-1)V^{*}(0)C\bar{u}^T(0).$$

Link must transform as

$$U_y(L-1) \to V(L-1)U_y(L-1)V^T(0)$$

- Link parallel to boundary on on other side  $(y \ge L)$  must then transform as:
- $U_x(x,y,..) \to V^*(x,y,..)U_x(x,y,..)V^T(x+1,y,..)$
- Gauge fields therefore obey complex-conjugate BCs.

#### The Two-Flavor Method



 Two fermion fields on each site indexed by flavor index:

$$\psi^{(1)}(x) = d(x), \ \psi^{(2)}(x) = C\bar{u}^T(x)$$

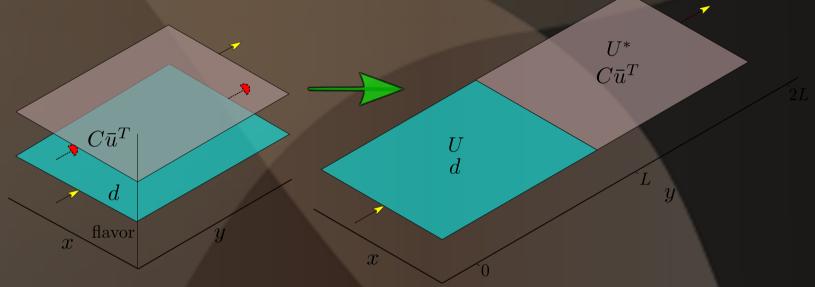
• BCs are:

$$\psi^{(1)}(x + L\hat{y}) = \psi^{(2)}(x),$$
  
$$\psi^{(2)}(x + L\hat{y}) = -\psi^{(1)}(x),$$

- Periodic BCs in other dirs.
- Single U-field shared by both flavors, with complex conj BCs.
- Dirac op for  $\psi^{(2)}$  uses  $U_{\mu}^{*}$ .

#### The One-Flavor Method

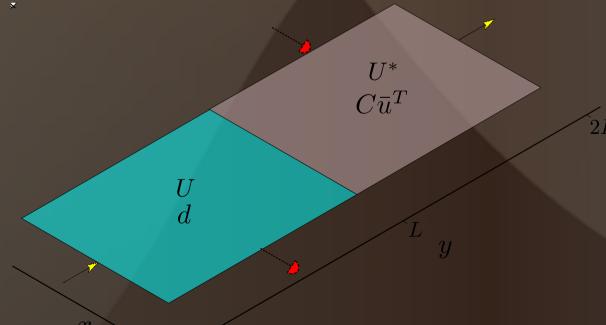
• Obtain equivalent formulation by unwrapping flavor indices onto two halves of doubled lattice:



- Antiperiodic boundary conditions in G-parity direction.
- U-field on first half and  $U^*$ -field on second half.

# Choosing an Approach

- One flavor setup is much easier to implement.
- However recall that we needed APBC in 2 directions for physical kinematics in  $\Delta I = 3/2$  calculation.
- G-parity in >1 dir using one-flavor method requires doubling the lattice again, which is highly inefficient.
- A second approach requires non-nearest neighbour communication:



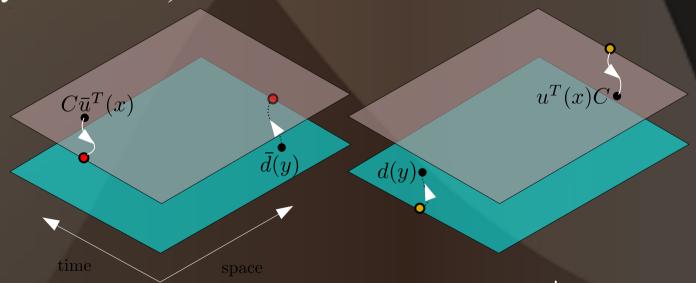
- Also inefficient depending on machine architecture.
- Choose to implement two-flavor method.

#### Unusual Contractions

 Flavor mixing at boundary allows contraction of up and down fields:

$$\psi_x^{(2)} \bar{\psi}_y^{(1)} = \mathcal{G}_{x,y}^{(2,1)} = C \bar{u}_x^T \bar{d}_y, 
\psi_y^{(1)} \bar{\psi}_x^{(2)} = \mathcal{G}_{y,x}^{(1,2)} = -\bar{d}_y u_x^T C^T$$

• Interpret as boundary creating/destroying flavor (violating baryon number):



• Also have  $\gamma^5$ -hermiticity:  $\left[\gamma^5 \mathcal{G}_{x,y}^{(2,1)} \gamma^5\right]^\dagger = \mathcal{G}_{y,x}^{(1,2)}$ 

# Exploiting the Underlying Gauge-Field Symmetry

- Quarks on flavor-1 plane interact with U field, and those on flavor-2 plane with U\*.
- Suggests propagators are related in some way.
- In fact, we find that:

$$\mathcal{G}_{x,z}^{(2,2)} = -\gamma^5 C \left[ \mathcal{G}_{x,z}^{(1,1)} \right]^* C \gamma^5$$

$$\mathcal{G}_{x,z}^{(1,2)} = +\gamma^5 C \left[ \mathcal{G}_{x,z}^{(2,1)} \right]^* C \gamma^5$$

- Relative sign due to sign at boundary between u and d.
- Substantially simplifies contractions.
- In some cases these relations can be used to reduce the number of propagator inversions required.

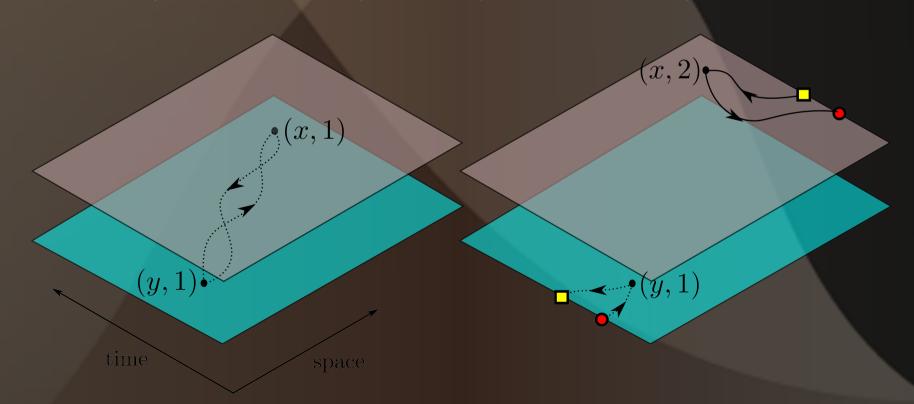
#### Pion Correlation Functions

•  $\pi^+$  correlation function

$$\langle \bar{d}_x \gamma^5 u_x \bar{u}_y \gamma^5 d_y \rangle = \langle \bar{\psi}_x^{(1)} [\gamma^5 C] \bar{\psi}_x^{(2)} {}^T \psi_y^{(2)} {}^T [C \gamma^5] \psi_y^{(1)} \rangle$$

• Now has *two* contractions:

$$\operatorname{tr}\left\{\mathcal{G}_{x,y}^{(1,1)\,\dagger}\mathcal{G}_{x,y}^{(1,1)}\right\} - \operatorname{tr}\left\{\mathcal{G}_{x,y}^{(2,1)\,\dagger}\mathcal{G}_{x,y}^{(2,1)}\right\}$$



# Locality

- Theory has one too many flavors. Must take square-root of  $s^\prime/s$  determinant in evolution to revert to 3 flavors.
- Determinant becomes non-local.
- Non-locality is however only a boundary effect that vanishes as  $L \to \infty$ . With sufficiently large volumes the effect should be minimal.
- Estimate size of effect?
  - Staggered ChPT?
  - Observe effect of changing from  $d \to C\bar{u}^T \to -d$  to  $d \to C\bar{u}^T \to +d$  for which  $\sqrt{\mathrm{Det}(D)}$  is local (= Pfaffian(D))?

# Charged Kaon Correlator

- $K^+$  analogue:  $|K^+'\rangle = \frac{1}{\sqrt{2}}(\bar{u}\gamma^5s \bar{s}'\gamma^5d)|0\rangle$
- 2-point function also has 4 contractions: (flavour indices 3 = s,  $4 = C\bar{s}'^T$ ):

$$\frac{1}{2} \operatorname{tr} \left\{ \mathcal{G}_{x,y}^{(3,3)} {}^{\dagger} \mathcal{G}_{x,y}^{(1,1)} \right\} + \frac{1}{2} \operatorname{tr} \left\{ \mathcal{G}_{x,y}^{(3,3)} \mathcal{G}_{x,y}^{(1,1)} {}^{\dagger} \right\}$$

$$+ \frac{1}{2} \operatorname{tr} \left\{ \mathcal{G}_{x,y}^{(4,3)} {}^{\dagger} \mathcal{G}_{x,y}^{(2,1)} \right\} + \frac{1}{2} \operatorname{tr} \left\{ \mathcal{G}_{x,y}^{(4,3)} \mathcal{G}_{x,y}^{(2,1)} {}^{\dagger} \right\}$$

- If we make the masses of the (s', s) and (u, d) doublets the same this is just the  $\pi^+$  correlation function but with the *opposite sign* between the contractions.
- Periodicity of spatial dependence appears to arise due to some cancellation between the two contractions.